

Lodge Pole Pine

REFERENCE MATERIAL

Pedigree

Institution: Tricon Lumber Received at INL: 2014

Location: Mineral County, MT Sample Preparation: Ground 2-inch chips to pass through a 1-inch sieve

Harvested: 2014 using a Vermeer BG480 grinder then material was dried in a drum dryer

Composition

Table 1. Chemical composition^a of Reference Lodge Pole Pine (mean of analyses completed 2/2015 & 4/2015)

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%Structural Ash	%Extractable Inorganics	%Water Extracted Glucan ^b	%Water Extracted Xylan ^b	%Water Extractives Others
0.27	0.13	0.13	0.04	2.45
%EtOH Extractives	%Lignin	%Glucan	%Xylan	%Galactan
1.74	30.50	41.37	5.90	2.84
%Arabinan+Mannan ^c	%Acetate	%Total		
12.20	1 17	00.00		

^aDetermined using NREL "Summative Mass Closure" LAP (NREL/TP-510-48087)

Proximate, Ultimate & Calorimetry

Table 2. Proximate, ultimate, and calorific values for Reference Lodge Pole Pine (reported on a dry basis; completed 3/2015)

Proximate ^a			Ultimate ^b			Calorimetry ^c	
%Volatile	%Ash	%Fixed Carbon	%Hydrogen	%Carbon	%Nitrogen	HHV	LHV
84.50	1.08	14.41	6.06	50.14	Below detection limit	8760	7371

^aProximate analysis was done according to ASTM D 5142-09

^bDetermined by HPLC following an acid hydrolysis of the water extractives

^c%Arabinan value includes %mannan, because arabinose and mannose co-elute on the HPLC column

^bUltimate analysis was conducted using a modified ASTM D5373-10 method (Flour and Plant Tissue Method) that uses a slightly different burn profile

^cHeating values (HHV, LHV) were determined with a calorimeter using ASTM D5865-10

Elemental Ash

Table 3. Elemental ash composition^a of Reference Lodge Pole Pine (completed 4/2015)

%Al as Al ₂ O ₃	%Ca as CaO	%Fe as Fe₂O₃	%K as K₂O		%Mn as MnO	%Na as Na₂O		%Si as SiO₂	%Ti as TiO₂	%S as SO₃
5.19	10.98	6.87	7.63	3.24	0.53	6.94	1.57	44.26	0.23	3.19

^aDetermined as described in ASTM standards D3174, D3682 and D6349

Lignin Chemistry

Table 4. Lignin chemistry of Reference Lodge Pole Pine (completed 1/2016)

Monolignol Composition ^a				Linkage /	Analysis ^b	
p-Hydroxyphenyl (H) content (% of total H+G+S)	Guaiacyl (G) content (% of total H+G+S)	Syringyl (S) content (% of total H+G+S)	ß-aryl ether (ß-O-4) (fraction of total) ^c	Phenylcoumaran (ß-5/a-O-4) (fraction of total)	Resinol (ß-ß) (fraction of total)	Dibenzodioxocin (5-5/4-O-ß) (fraction of total)
0	100	0	73	21	4	1

^aDetermined by integration of peak volumes of ball-milled whole cell wall samples, swelled in 4:1 DMSO:Py, and analyzed by gel-state HSQC NMR (Mansfield, S. D., et al. (2012) Nature Protocols, 7(9), 1579-1589)

^aDetermined as described in ASTM standards D3174, D3682 and D6349

^bQuantitative data on the different types of chemical linkages between monolignols in a biomass sample. Determined by integrating peak volumes in solution-state HSQC NMR spectra of acetylated whole cell wall samples

^cEther bond between the ß carbon on one monolignol to the phenolic oxygen on a second monolignol. This is typically the most common linkage found in native lignin samples (Vanholme, R., et al. (2010) Plant Physiol., 153, 895-905)

Particle Characteristics

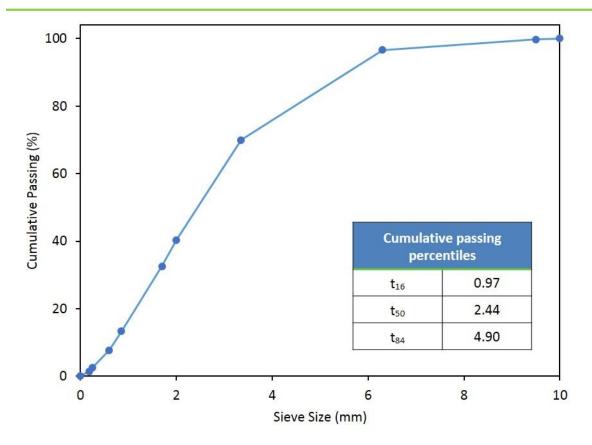


Figure 1. Cumulative passing percent of 1-inch Reference Lodge Pole Pine determined according to ANSI/ASAE S319.4 using a Ro-Tap test sieve shaker (Model RX-29, W.S. Tyler) and a 15 minute total sieving time (completed 4/2015). The cumulative passing percentile sieve sizes (e.g., t_{16}) were calculated by interpolation and represent theoretical sieve sizes that would retain 16, 50 or 84% of the particles by mass.

Contact

For questions regarding biomass material or analytical data please contact Amber Hoover at amber.hoover@inl.gov or 208-526-5992.

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